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***** Welcome to STN International *****

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
NEWS 2 "Ask CAS" for self-help around the clock
NEWS 3 SEP 01 New pricing for the Save Answers for SciFinder Wizard within

NEWS 4 OCT 28 STN Express with Discover!
NEWS 5 NOV 30 KOREAPAT now available on STN
NEWS 6 DEC 01 PHAR reloaded with additional data
NEWS 7 DEC 01 LISA now available on STN
NEWS 7 DEC 09 12 databases to be removed from STN on December 31, 2004

NEWS 8 DEC 15 MEDLINE update schedule for December 2004
NEWS 9 DEC 17 ELCOM reloaded; updating to resume; current-awareness

NEWS 10 DEC 17 alerts (SDIs) affected
COMPUAB reloaded; updating to resume; current-awareness

NEWS 11 DEC 17 alerts (SDIs) affected
SOLIDSTATE reloaded; updating to resume; current-awareness

NEWS 12 DEC 17 alerts (SDIs) affected
CERAB reloaded; updating to resume; current-awareness

NEWS 13 DEC 17 alerts (SDIs) affected
THREE NEW FIELDS ADDED TO IFIPAT/IFIUDB/IFICDB

NEWS EXPRESS OCTOBER 29 CURRENT WINDOWS VERSION IS V7.01A, CURRENT
MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
AND CURRENT DISCOVER FILE IS DATED 11 AUGUST 2004

NEWS HOURS STN Operating Hours Plus Help Desk Availability
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NEWS PHONE Direct Dial and Telecommunication Network Access to STN
NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

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***** STN Columbus *****

FILE 'HOME' ENTERED AT 10:25:10 ON 28 DEC 2004

=> FIL REGISTRY
COST IN U.S. DOLLARS
TOTAL
SINCE FILE
ENTRY
SESSION
FULL ESTIMATED COST 0.21

FILE 'REGISTRY' ENTERED AT 10:25:22 ON 28 DEC 2004
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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 26 DEC 2004 HIGHEST RN 802853-20-9
DICTIONARY FILE UPDATES: 26 DEC 2004 HIGHEST RN 802853-20-9

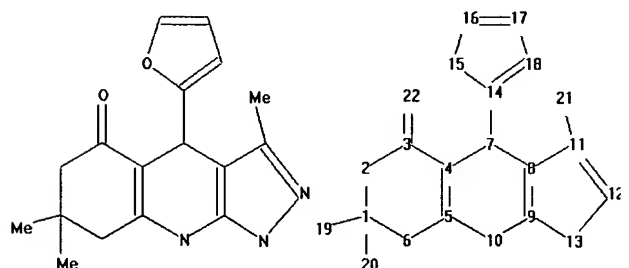
TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>
Uploading H:\STN queries\10612885b.str



chain nodes :
19 20 21 22
ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18
chain bonds :
1-19 1-20 3-22 7-14 11-21
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-10 7-8 8-9 8-11 9-10 9-13
11-12 12-13 14-15 14-18 15-16 16-17 17-18
exact/norm bonds :
1-2 1-6 2-3 3-4 3-22 4-5 4-7 5-6 5-10 7-8 8-9 8-11 9-10
9-13 11-12 12-13 14-15 14-18 15-16 16-17 17-18
exact bonds :
1-19 1-20 7-14 11-21

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom
9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom
17:Atom 18:Atom 19:CLASS 20:CLASS 21:CLASS 22:CLASS

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR

Page 1-A



Page 2-A

NODE ATTRIBUTES:
HCOUNT IS M3 AT 19
HCOUNT IS M3 AT 20
HCOUNT IS M3 AT 21
NSPEC IS R AT 1
NSPEC IS R AT 2
NSPEC IS R AT 3
NSPEC IS R AT 4
NSPEC IS R AT 5
NSPEC IS R AT 6
NSPEC IS R AT 7
NSPEC IS R AT 8
NSPEC IS R AT 9
NSPEC IS R AT 10
NSPEC IS R AT 11
NSPEC IS R AT 12
NSPEC IS R AT 13
NSPEC IS R AT 14
NSPEC IS R AT 15
NSPEC IS R AT 16
NSPEC IS R AT 17
NSPEC IS R AT 18
NSPEC IS C AT 19
NSPEC IS C AT 20

NSPEC IS C AT 21
NSPEC IS C AT 22
DEFAULT MLEVEL IS ATOM
MLEVEL IS CLASS AT 19 20 21 22
DEFAULT ELEVEL IS LIMITED

GRAPH ATTRIBUTES:
RING(S) ARE ISOLATED OR EMBEDDED
NUMBER OF NODES IS 22

STEREO ATTRIBUTES: NONE

=> s l1 sam
SAMPLE SEARCH INITIATED 10:25:56 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0
ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 0 TO 0
PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s l1 fam
SAMPLE SEARCH INITIATED 10:26:03 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0
ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 0 TO 0
PROJECTED ANSWERS: 0 TO 0

L3 0 SEA FAM SAM L1

=> s l1 fu1
FULL SEARCH INITIATED 10:26:08 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 21 TO ITERATE

100.0% PROCESSED 21 ITERATIONS 8
ANSWERS
SEARCH TIME: 00.00.01

L4 8 SEA SSS FUL L1

=> s l1 fu1 fam
FULL SEARCH INITIATED 10:26:12 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 5 TO ITERATE

100.0% PROCESSED 5 ITERATIONS 1
ANSWERS
SEARCH TIME: 00.00.01

L5 1 SEA FAM FUL L1

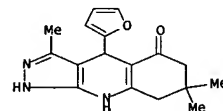
=> l5 not l4
L6 0 L5 NOT L4

=> d l5 1 a11

L5 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2004 ACS on STN
RN 645337-25-3 REGISTRY
ED Entered STN: 03 Feb 2004
CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl- (9CI) (CA INDEX NAME)
FS 3D CONCORD
MF C17 H19 N3 O2
SR CA
LC STN Files: CA, CAPLUS, USPATFULL
DT,CA Caplus document type: Patent
RL,P Roles from patents: BIOL (Biological study); USES (Uses)

Ring System Data

Elemental Analysis EA	Elemental Sequence ES	Size of the Rings SZ	Ring System Formula RF	Ring Identifier RID	RID Occurrence Count
C40	10C4	5	C40	16.138.5	1
C3N2-C5N-C6	N2C3-NC5-C6	5-6-6	C10N3	1894.85.24	1



Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	1	pH 1	(1) ACD
Bioconc. Factor (BCF)	8.39	pH 4	(1) ACD
Bioconc. Factor (BCF)	97.8	pH 7	(1) ACD

Bioconc. Factor (BCF) 98.8 pH 8 (1) ACD
Bioconc. Factor (BCF) 98.8 pH 10 (1) ACD
Boiling Point (BP) 478.2+/-45.0 deg C 760 Torr (1) ACD
Enthalpy of Vap. (HVAP) 74.23+/-3.0 kJ/mol (1) ACD
Flash Point (FP) 243.0+/-51.7 deg C (1) ACD
Freely Rotatable Bonds (FRB) 1 (1) ACD
H acceptors (HAC) 5 (1) ACD
H donors (HD) 2 (1) ACD
Koc (KOC) 1 pH 1 (1) ACD
Koc (KOC) 79.0 pH 4 (1) ACD
Koc (KOC) 922 pH 7 (1) ACD
Koc (KOC) 931 pH 8 (1) ACD
Koc (KOC) 932 pH 10 (1) ACD
logD (LOGD) -0.49 pH 1 (1) ACD
logD (LOGD) 1.86 pH 4 (1) ACD
logD (LOGD) 2.92 pH 7 (1) ACD
logD (LOGD) 2.93 pH 8 (1) ACD
logD (LOGD) 2.93 pH 10 (1) ACD
logP (LOGP) 2.928+/-0.412 (1) ACD
Molar Solubility (SLB.MOL) >=0.1 - <1 mol/L pH 1 (1) ACD
Molar Solubility (SLB.MOL) <0.01 mol/L pH 4 (1) ACD
Molar Solubility (SLB.MOL) <0.01 mol/L pH 7 (1) ACD
Molar Solubility (SLB.MOL) <0.01 mol/L pH 8 (1) ACD
Molar Solubility (SLB.MOL) <0.01 mol/L pH 10 (1) ACD
Molecular Weight (MW) 297.35 (1) ACD
pKa (PKA) 13.94+/-0.60 Most Acidic (1) ACD
pKa (PKA) 4.82+/-0.60 Most Basic (1) ACD
Vapor Pressure (VP) 2.62E-09 Torr 25 deg C (1) ACD

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software
Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1

AN 140:87744 CA Full-text
TI Affinity small molecules for the EPO receptor
IN Olsson, Lennart; Naranda, Tatjana
PA Recepton, Inc., USA
SO PCT Int. Appl., 85 pp.
CODEN: PIXX02
DT Patent
LA English
IC ICM C07K
CC 1-12 (Pharmacology)
Section cross-reference(s): 2

FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004005323	A2	20040115	WO 2003-US21394	20030703
WO 2004005323	A3	20040701		

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZM, ZW
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG
US 2004171541 A1 20040902 US 2003-613754 20030702
US 2004116346 A1 20040617 US 2003-612885 20030703
PRAI US 2002-393360P 20020703
US 2002-393361P 20020703
US 2002-394110P 20020703
AB Compds. are provided that complex with the modulating domain of erythropoietin receptor (EPO-R) for use with EPO-R to determine the presence of EPO-R, the ability of other mols. to bind to the modulating domain in competitive assays and to induce a signal by EPO-R into a cell when bound by the subject compds. in a physiol. environment. The compds. are characterized by having a six-membered heterocyclic ring comprising at least one nitrogen atom and include substituted triazopyrimidine, pyridazinone, pyridine and piperidine.
ST EPO receptor modulator small mol
IT Proteins
RL: BSU (Biological study, unclassified); BIOL (Biological study) (Bcl-XL, expression; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
IT Peptides, biological studies
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(EPO receptor modulating sequence; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
IT Cell membrane
(EPO receptors of; affinity small mols. for erythropoietin

(EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Anemia (disease)
Cell proliferation
Combinatorial library
Drug delivery systems
Drug screening
Erythrocyte
Erythropoiesis
Hematocrit
Hematopoietic precursor cell
Human
Reticulocyte
(affinity small mols. for erythropoietin (EPO) receptor and

EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Erythropoietin receptors
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and

EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Nerve
(neuron; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Cytoprotective agents
(neuroprotective; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT 2503-56-2 40775-78-8 51646-16-3 51646-17-4 51646-19-6 51646-43-6
56347-20-7 63901-48-4 90559-98-1 90815-61-5 113967-71-8 113967-74-1 194342-06-8 212074-47-0 244167-89-3 245082-87-5 245413-82-5 259683-29-9 261704-08-9 261704-09-0 262291-81-6

263267-38-5 287728-46-5 303145-64-4 303145-73-5 338793-16-1 645337-19-5 645337-20-8 645337-21-9 645337-22-0 645337-23-1 645337-24-2 645337-25-3
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and

EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT 11096-26-7 Erythropoietin
RL: BSU (Biological study, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and

EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT 239133-03-0 645415-22-1
RL: PRP (Properties)
(unclaimed sequence; affinity small mols. for the EPO receptor)

=> d 14 1-8 ful
'FUL' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN
SAM - Index Name, MF, and structure - no RN
FIDE - All substance data, except sequence data
IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data
SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used
SQD - Protein sequence data, includes RN
SQD3 - Same as SQD, but 3-letter amino acid codes are used
SQN - Protein sequence name information, includes RN

CALC - Table of calculated properties
EPROP - Table of experimental properties
PROP - EPROP and CALC

Any CA File format may be combined with any substance format to

obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract
APPS -- Application and Priority Information
BIB -- CA Accession Number, plus Bibliographic Data
CAN -- CA Accession Number
CBIB -- CA Accession Number, plus Bibliographic Data (compressed)
IND -- Index Data
IPC -- International Patent Classification
PATS -- PI, SO
STD -- BIB, IPC, and NCL

IABS -- ABS, indented, with text labels
IBIB -- BIB, indented, with text labels
ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)
OIBIB ----- OBIB, indented with text labels

SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.
The MAX format is the same as ALL.
The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

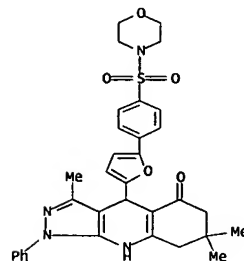
For additional information, please consult the following help messages:

HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):all

L4 ANSWER 1 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
RN 748146-78-3 REGISTRY
ED Entered STN: 20 Sep 2004
CN INDEX NAME NOT YET ASSIGNED
FS 3D CONCORD
MF C33 H34 N4 O5 S
SR Chemical Library

Ring System Data

Elemental Analysis EA	Elemental Sequence ES	Size of the Rings SZ	Ring System Formula RF	Ring Identifier RID	RID Occurrence Count
C4O	OC4	5	C4O	16.138.5	1
C6	C6	6	C6	46.150.18	2
C4NO	NC2OC2	6	C4NO	46.402.1	1
C3N2-C5N-C6	N2C3-NC5-C6	5-6-6	C10N3	1894.85.24	1



Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	12.13	pH 1	(1) ACD
Bioconc. Factor (BCF)	649	pH 4	(1) ACD
Bioconc. Factor (BCF)	1353	pH 7	(1) ACD
Bioconc. Factor (BCF)	1354	pH 8	(1) ACD
Bioconc. Factor (BCF)	1355	pH 10	(1) ACD
Boiling Point (BP)	749.7+/-70.0 deg C	760 Torr	(1) ACD
Enthalpy of Vap. (HVAP)	109.28+/-3.0 kJ/mol		(1) ACD
Flash Point (FP)	407.2+/-64.2 deg C		(1) ACD
Freely Rotatable Bonds (FRB)	4		(1) ACD
H acceptors (HAC)	9		(1) ACD
H donors (HD)	1		(1) ACD
Koc (KOC)	9.55	pH 1	(1) ACD
Koc (KOC)	2910	pH 4	(1) ACD
Koc (KOC)	6062	pH 7	(1) ACD
Koc (KOC)	6067	pH 8	(1) ACD
Koc (KOC)	6068	pH 10	(1) ACD
logD (LOGD)	1.62	pH 1	(1) ACD
logD (LOGD)	4.10	pH 4	(1) ACD
logD (LOGD)	4.42	pH 7	(1) ACD
logD (LOGD)	4.42	pH 8	(1) ACD
logD (LOGD)	4.42	pH 10	(1) ACD
logP (LOGP)	4.424+/-0.655		(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1) ACD

Molar Solubility (SLB.MOL) <0.01 mol/L pH 10 (1) ACD
 Molecular Weight (MW) 598.71 (1) ACD
 pKa (PKA) 3.95+/-0.60 Most Basic (1) ACD
 Vapor Pressure (VP) 2.33E-22 Torr 25 deg C (1) ACD

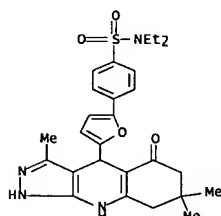
(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software
 Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

L4 ANSWER 2 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 748146-41-0 REGISTRY
 ED Entered STN: 20 Sep 2004
 CN INDEX NAME NOT YET ASSIGNED
 FS 3D CONCORD
 MF C27 H32 N4 O4 S
 SR Chemical Library

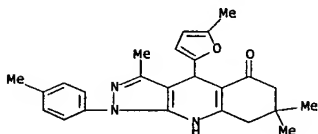
Ring System Data

Elemental Analysis EA	Elemental Sequence ES	Size of the Rings SZ	Ring System Formula RF	Ring Identifier RID	RID Occurrence Count
C40	OC4	5	C40	16.138.5	1
C6	C6	6	C6	46.150.18	1
C3N2-C5N-C6	N2C3-NC5-C6	5-6-6	C10N3	1894.85.24	1



Calculated Properties (CALC)

C40	OC4	5	C40	16.138.5	1
C6	C6	6	C6	46.150.18	1
C3N2-C5N-C6	N2C3-NC5-C6	5-6-6	C10N3	1894.85.24	1



Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	2.22	pH 1	(1) ACD
Bioconc. Factor (BCF)	274	pH 4	(1) ACD
Bioconc. Factor (BCF)	3190	pH 7	(1) ACD
Bioconc. Factor (BCF)	3221	pH 8	(1) ACD
Bioconc. Factor (BCF)	3225	pH 10	(1) ACD
Boiling Point (BP)	680.7+/-65.0 deg C	760 Torr	(1) ACD
Enthalpy of Vap. (HVAP)	99.88+/-3.0 kJ/mol		(1) ACD
Flash Point (FP)	365.5+/-61.7 deg C		(1) ACD
Freely Rotatable Bonds (FRB)	5		(1) ACD
H acceptors (HAC)	8		(1) ACD
H donors (HD)	2		(1) ACD
Koc (KOC)	4.28	pH 1	(1) ACD
Koc (KOC)	958	pH 4	(1) ACD
Koc (KOC)	11170	pH 7	(1) ACD
Koc (KOC)	11278	pH 8	(1) ACD
Koc (KOC)	11289	pH 10	(1) ACD
logD (LOGD)	1.50	pH 1	(1) ACD
logD (LOGD)	3.85	pH 4	(1) ACD
logD (LOGD)	4.91	pH 7	(1) ACD
logD (LOGD)	4.92	pH 8	(1) ACD
logD (LOGD)	4.92	pH 10	(1) ACD
logP (LOGP)	4.920+/-0.532		(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 10	(1) ACD
Molecular Weight (MW)	598.63		(1) ACD
pKa (PKA)	13.94+/-0.60	Most Acidic	(1) ACD
pKa (PKA)	4.82+/-0.60	Most Basic	(1) ACD
Vapor Pressure (VP)	2.19E-18 Torr	25 deg C	(1) ACD

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	1.22	pH 1	(1) ACD
Bioconc. Factor (BCF)	274	pH 4	(1) ACD
Bioconc. Factor (BCF)	3190	pH 7	(1) ACD
Bioconc. Factor (BCF)	3221	pH 8	(1) ACD
Bioconc. Factor (BCF)	3225	pH 10	(1) ACD
Boiling Point (BP)	680.7+/-65.0 deg C	760 Torr	(1) ACD
Enthalpy of Vap. (HVAP)	99.88+/-3.0 kJ/mol		(1) ACD
Flash Point (FP)	365.5+/-61.7 deg C		(1) ACD
Freely Rotatable Bonds (FRB)	5		(1) ACD
H acceptors (HAC)	8		(1) ACD
H donors (HD)	2		(1) ACD
Koc (KOC)	4.28	pH 1	(1) ACD
Koc (KOC)	958	pH 4	(1) ACD
Koc (KOC)	11170	pH 7	(1) ACD
Koc (KOC)	11278	pH 8	(1) ACD
Koc (KOC)	11289	pH 10	(1) ACD
logD (LOGD)	1.50	pH 1	(1) ACD
logD (LOGD)	3.85	pH 4	(1) ACD
logD (LOGD)	4.91	pH 7	(1) ACD
logD (LOGD)	4.92	pH 8	(1) ACD
logD (LOGD)	4.92	pH 10	(1) ACD
logP (LOGP)	4.920+/-0.532		(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 10	(1) ACD
Molecular Weight (MW)	598.63		(1) ACD
pKa (PKA)	13.94+/-0.60	Most Acidic	(1) ACD
pKa (PKA)	4.82+/-0.60	Most Basic	(1) ACD
Vapor Pressure (VP)	2.19E-18 Torr	25 deg C	(1) ACD

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software
 Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

L4 ANSWER 3 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 748145-15-5 REGISTRY
 ED Entered STN: 20 Sep 2004
 CN INDEX NAME NOT YET ASSIGNED
 FS 3D CONCORD
 MF C25 H27 N3 O2
 SR Chemical Library

Ring System Data

Elemental Analysis EA	Elemental Sequence ES	Size of the Rings SZ	Ring System Formula RF	Ring Identifier RID	RID Occurrence Count
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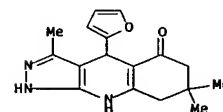
(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software
 Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

L4 ANSWER 4 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 645337-25-3 REGISTRY
 ED Entered STN: 03 Feb 2004
 CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl- (9CI) (CA INDEX NAME)
 FS 3D CONCORD
 MF C17 H19 N3 O2
 SR CA
 LC STN Files: CA, CAPLUS, USPATFULL
 DT.CA Caplus document type: Patent
 RL.P Roles from patents: BIOL (Biological study); USES (Uses)

Ring System Data

Elemental Analysis EA	Elemental Sequence ES	Size of the Rings SZ	Ring System Formula RF	Ring Identifier RID	RID Occurrence Count
C40	OC4	5	C40	16.138.5	1
C3N2-C5N-C6	N2C3-NC5-C6	5-6-6	C10N3	1894.85.24	1



Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	1	pH 1	(1) ACD
Bioconc. Factor (BCF)	18.39	pH 4	(1) ACD
Bioconc. Factor (BCF)	97.8	pH 7	(1) ACD
Bioconc. Factor (BCF)	98.8	pH 8	(1) ACD
Bioconc. Factor (BCF)	98.8	pH 10	(1) ACD
Boiling Point (BP)	478.2+/-45.0 deg C	760 Torr	(1) ACD

Enthalpy of Vap. (HVAP) 74.23+/-3.0 kJ/mol (1) ACD
Flash Point (FP) 243.0+/-51.7 deg C (1) ACD
Freely Rotatable Bonds (FRB) 1 (1) ACD
H acceptors (HAC) 5 (1) ACD
H donors (HD) 2 (1) ACD
Koc (KOC) 1 pH 1 (1) ACD
Koc (KOC) 79.0 pH 4 (1) ACD
Koc (KOC) 922 pH 7 (1) ACD
Koc (KOC) 931 pH 8 (1) ACD
Koc (KOC) 932 pH 10 (1) ACD
logD (LOGD) -0.49 pH 1 (1) ACD
logD (LOGD) 1.86 pH 4 (1) ACD
logD (LOGD) 2.92 pH 7 (1) ACD
logD (LOGD) 2.93 pH 8 (1) ACD
logD (LOGD) 2.93 pH 10 (1) ACD
logP (LOGP) 2.928+/-0.412 (1) ACD
Molar Solubility (SLB.MOL) >=0.1 - <1 mol/L pH 1 (1) ACD
Molar Solubility (SLB.MOL) <0.01 mol/L pH 4 (1) ACD
Molar Solubility (SLB.MOL) <0.01 mol/L pH 7 (1) ACD
Molar Solubility (SLB.MOL) <0.01 mol/L pH 8 (1) ACD
Molar Solubility (SLB.MOL) <0.01 mol/L pH 10 (1) ACD
Molecular Weight (MW) 297.35 (1) ACD
pKa (PKA) 13.94+/-0.60 Most Acidic (1) ACD
pKa (PKA) 4.82+/-0.60 Most Basic (1) ACD
Vapor Pressure (VP) 2.62E-09 Torr 25 deg C (1) ACD

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software
Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1

AN 140:87744 CA Full-text
TI Affinity small molecules for the EPO receptor
IN Olsson, Lennart; Naranda, Tatjana
PA Receptor, Inc., USA
SO PCT Int. Appl., 85 pp.
CODEN: PIXX02
DT Patent
LA English
IC ICM C07K
CC 1-12 (Pharmacology)
Section cross-reference(s): 2
FAN.CNT 1
PATENT NO. KIND DATE APPLICATION NO. DATE
PI WO 2004005323 A2 20040115 WO 2003-US21394 20030703
WO 2004005323 A3 20040701
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD,

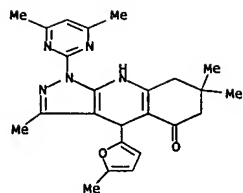
the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
IT Anemia (disease)
Cell proliferation
Combinatorial library
Drug delivery systems
Drug screening
Erythrocyte
Erythropoiesis
Hematocrit
Hematopoietic precursor cell
Human
Reticulocyte
(affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
IT Erythropoietin receptors
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
IT Nerve
(neuron; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
IT Cytoprotective agents
(neuroprotective; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
IT 2503-56-2 40775-78-8 51646-16-3 51646-17-4 51646-19-6 51646-43-6
56347-20-7 63901-48-4 90559-98-1 90815-61-5 113967-71-8 113967-74-1 194342-06-8 212074-47-0 244167-89-3 245082-87-5 245413-82-5 259683-29-9 261704-08-9 261704-09-0 262291-81-6 263267-38-5 287728-46-5 303145-64-4 303145-73-5 338793-16-1 645337-19-5 645337-20-8 645337-21-9 645337-22-0 645337-

GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZM, ZW
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG
US 2004171541 A1 20040902 US 2003-613754 20030702
US 2004116346 A1 20040617 US 2003-612885 20030703
PRAI US 2002-393360P 20020703
US 2002-393361P 20020703
US 2002-394110P 20020703
AB Comps. are provided that complex with the modulating domain of erythropoietin receptor (EPO-R) for use with EPO-R to determine the presence of EPO-R, the ability of other mols. to bind to the modulating domain in competitive assays and to induce a signal by EPO-R into a cell when bound by the subject compds. in a physiol. environment. The compds. are characterized by having a six-membered heterocyclic ring comprising at least one nitrogen atom and include substituted triazolopyrimidine, pyridazinone, pyridine and piperidine.
ST EPO receptor modulator small mol
IT Proteins
RL: BSU (Biological study, unclassified); BIOL (Biological study) (Bcl-xL, expression; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
IT Peptides, biological studies
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(EPO receptor modulating sequence; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
IT Cell membrane
(EPO receptors of; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating

23-1 645337-24-2 645337-25-3
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
IT 11096-26-7, Erythropoietin
RL: BSU (Biological study, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
IT 239133-03-0 645415-22-1
RL: PRP (Properties)
(unclaimed sequence; affinity small mols. for the EPO receptor)
L4 ANSWER 5 OF 8 REGISTRY COPYRIGHT 2004 ACS ON STN
RN 521318-71-8 REGISTRY
ED Entered STN: 28 May 2003
CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 1-(4,6-dimethyl-2-pyrimidinyl)-
1,4,6,7,8,9-hexahydro-3,7,7-trimethyl-4-(5-methyl-2-furanyl)-(9CI)
(CA
INDEX NAME)
FS 3D CONCORD
MF C24 H27 N5 O2
SR Chemical Library
LC STN Files: CHEMCATS

Ring System Data

Elemental Analysis	Elemental Sequence	Size of the Rings	Ring System Formula	Ring Identifier	RID Occurrence Count
EA	ES	SZ	RF	RID	
C40	OC4	5	C40	16.138.5	1
C4N2	ICNC3	6	C4N2	46.195.39	1
C3N2-C5N-C6	N2C3-NC5-C6	5-6-6	C10N3	1894.85.24	1



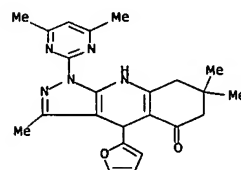
Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

L4 ANSWER 6 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 521284-01-5 REGISTRY
 ED Entered STN: 28 May 2003
 CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 1-(4,6-dimethyl-2-pyrimidinyl)-4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl- (9CI) (CA INDEX NAME)
 FS 3D CONCORD
 MF C23 H25 N5 O2
 SR Chemical Library
 LC STN Files: CHEMCATS

Ring System Data

Elemental Analysis EA	Elemental Sequence ES	Size of the Rings SZ	Ring System Formula RF	Ring Identifier RID	RID Occurrence Count
C40	OC4	5	C40	16.138.5	1
C4N2	NCNC3	6	C4N2	46.195.39	1
C3N2-C5N-C6	N2C3-NC5-C6	5-6-6	C10N3	1894.85.24	1



Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	108	pH 1	(1) ACD
Bioconc. Factor (BCF)	283	pH 4	(1) ACD
Bioconc. Factor (BCF)	283	pH 7	(1) ACD

Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	242	pH 1	(1) ACD
Bioconc. Factor (BCF)	633	pH 4	(1) ACD
Bioconc. Factor (BCF)	634	pH 7	(1) ACD
Bioconc. Factor (BCF)	634	pH 8	(1) ACD
Bioconc. Factor (BCF)	634	pH 10	(1) ACD
Boiling Point (BP)	607.6+/-65.0 deg C	760.0 Torr	(1) ACD
Enthalpy of Vap. (HVPAP)	90.29+/-3.0 kJ/mol		(1) ACD
Flash Point (FP)	321.2+/-61.7 deg C		(1) ACD
Freely Rotatable Bonds (FRB)	2		(1) ACD
H acceptors (HAC)	7		(1) ACD
H donors (HD)	1		(1) ACD
Koc (KOC)	1346	pH 1	(1) ACD
Koc (KOC)	3518	pH 4	(1) ACD
Koc (KOC)	3524	pH 7	(1) ACD
Koc (KOC)	3524	pH 8	(1) ACD
Koc (KOC)	3524	pH 10	(1) ACD
logD (LOGD)	3.57	pH 1	(1) ACD
logD (LOGD)	3.99	pH 4	(1) ACD
logD (LOGD)	3.99	pH 7	(1) ACD
logD (LOGD)	3.99	pH 8	(1) ACD
logD (LOGD)	3.99	pH 10	(1) ACD
logP (LOGP)	3.990+/-0.891		(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 10	(1) ACD
Molecular Weight (MW)	417.50		(1) ACD
pKa (PKA)	1.15+/-0.60	Most Basic	(1) ACD
Vapor Pressure (VP)	1.04E-14 Torr	25.0 deg C	(1) ACD

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software

Bioconc. Factor (BCF)	283	pH 8	(1) ACD
Bioconc. Factor (BCF)	283	pH 10	(1) ACD
Boiling Point (BP)	601.5+/-65.0 deg C	760.0 Torr	(1) ACD
Enthalpy of Vap. (HVPAP)	89.51+/-3.0 kJ/mol		(1) ACD
Flash Point (FP)	317.6+/-61.7 deg C		(1) ACD
Freely Rotatable Bonds (FRB)	2		(1) ACD
H acceptors (HAC)	7		(1) ACD
H donors (HD)	1		(1) ACD
Koc (KOC)	756	pH 1	(1) ACD
Koc (KOC)	1977	pH 4	(1) ACD
Koc (KOC)	1980	pH 7	(1) ACD
Koc (KOC)	1980	pH 8	(1) ACD
Koc (KOC)	1980	pH 10	(1) ACD
logD (LOGD)	3.11	pH 1	(1) ACD
logD (LOGD)	3.53	pH 4	(1) ACD
logD (LOGD)	3.53	pH 7	(1) ACD
logD (LOGD)	3.53	pH 8	(1) ACD
logD (LOGD)	3.53	pH 10	(1) ACD
logP (LOGP)	3.530+/-0.890		(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 10	(1) ACD
Molecular Weight (MW)	403.48		(1) ACD
pKa (PKA)	1.15+/-0.60	Most Basic	(1) ACD
Vapor Pressure (VP)	2.00E-14 Torr	25.0 deg C	(1) ACD

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software

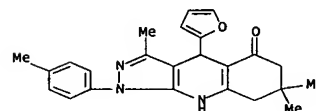
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See HELP PROPERTIES for information about property data sources in REGISTRY.

L4 ANSWER 7 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 380450-98-6 REGISTRY
 ED Entered STN: 04 Jan 2002
 CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl-1-(4-methylphenyl)- (9CI) (CA INDEX NAME)
 FS 3D CONCORD
 MF C24 H25 N3 O2
 SR Chemical Library
 LC STN Files: CHEMCATS

Ring System Data

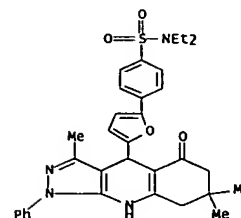
Elemental Analysis EA	Elemental Sequence ES	Size of the Rings SZ	Ring System Formula RF	Ring Identifier RID	RID Occurrence Count
C40	OC4	5	C40	16.138.5	1
C6	OC6	6	C6	46.150.18	1
C3N2-C5N-C6	N2C3-NC5-C6	5-6-6	C10N3	1894.85.24	1



L4 ANSWER 8 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 378189-53-8 REGISTRY
 ED Entered STN: 26 Dec 2001
 CN Benzenesulfonamide, N,N-diethyl-4-[5-(4,5,6,7,8,9-hexahydro-3,7,7-trimethyl-5-oxo-1-phenyl-1H-pyrazolo[3,4-b]quinolin-4-yl)-2-furanyl]- (9CI) (CA INDEX NAME)
 FS 3D CONCORD
 MF C33 H36 N4 O4 S
 SR Chemical Library
 LC STN Files: CHEMCATS

Ring System Data

Elemental Analysis EA	Elemental Sequence ES	Size of the Rings SZ	Ring System Formula RF	Ring Identifier RID	RID Occurrence Count
C40	OC4	5	C40	16.138.5	1
C6	OC6	6	C6	46.150.18	2
C3N2-C5N-C6	N2C3-NC5-C6	5-6-6	C10N3	1894.85.24	1



Calculated Properties (CALC)

PROPERTY (CODE)	VALUE	CONDITION	NOTE
Bioconc. Factor (BCF)	22.5	pH 1	(1) ACD
Bioconc. Factor (BCF)	6852	pH 4	(1) ACD
Bioconc. Factor (BCF)	14276	pH 7	(1) ACD
Bioconc. Factor (BCF)	14290	pH 8	(1) ACD
Bioconc. Factor (BCF)	14291	pH 10	(1) ACD
Boiling Point (BP)	712.3+/-70.0 deg C	760.0 Torr	(1) ACD
Enthalpy of Vap. (HVAP)	104.14+/-3.0 kJ/mol		(1) ACD
Flash Point (FP)	384.6+/-64.2 deg C		(1) ACD
Freely Rotatable Bonds (FRB)	6		(1) ACD
H acceptors (HAC)	8		(1) ACD
H donors (HD)	1		(1) ACD
Koc (KOC)	51.6	pH 1	(1) ACD
Koc (KOC)	15714	pH 4	(1) ACD
Koc (KOC)	32737	pH 7	(1) ACD
Koc (KOC)	32769	pH 8	(1) ACD
Koc (KOC)	32772	pH 10	(1) ACD
logD (LOGD)	2.97	pH 1	(1) ACD
logD (LOGD)	5.45	pH 4	(1) ACD
logD (LOGD)	5.77	pH 7	(1) ACD
logD (LOGD)	5.77	pH 8	(1) ACD
logD (LOGD)	5.77	pH 10	(1) ACD
logP (LOGP)	5.771+/-0.615		(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 1	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 4	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 7	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 8	(1) ACD
Molar Solubility (SLB.MOL)	<0.01 mol/L	pH 10	(1) ACD
Molecular Weight (MW)	584.73		(1) ACD
pKa (PKA)	3.95+/-0.60	Most Basic	(1) ACD
Vapor Pressure (VP)	3.87E-20 Torr	25.0 deg C	(1) ACD

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software
Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

=> d 14 1-8 ed

L4 ANSWER 1 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 20 Sep 2004

L4 ANSWER 2 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 20 Sep 2004

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=> d his

(FILE 'HOME' ENTERED AT 10:25:10 ON 28 DEC 2004)

FILE 'REGISTRY' ENTERED AT 10:25:22 ON 28 DEC 2004
STRUCTURE UPLOADED
L1 0 S L1 SAM
L2 0 S L1 FAM
L3 8 S L1 FUL
L4 1 S L1 FUL FAM
L5 0 L5 NOT L4
L6

FILE 'HCAPLUS, USPATFULL, CHEMCATS' ENTERED AT 10:35:54 ON 28 DEC 2004

=> s 14

L7 9 L4

=> d 19 ed

L9 NOT FOUND

The L-number entered has not been defined in this session, or it has been deleted. To see the L-numbers currently defined in this session, enter DISPLAY HISTORY at an arrow prompt (=>).

=> d 17 1-9 ed

NO VALID FORMATS ENTERED FOR FILE 'USPATFULL'
NO VALID FORMATS ENTERED FOR FILE 'CHEMCATS'
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REENTER DISPLAY FORMAT FOR ALL FILES (FILEDEFAULT):end

=> d 17 1-9 ibib abs

NO VALID FORMATS ENTERED FOR FILE 'CHEMCATS'
In a multiframe environment, each file must have at least one valid format requested. Refer to file specific help messages or the STNGUIDE file for information on formats available in individual files.

REENTER DISPLAY FORMAT FOR ALL FILES (FILEDEFAULT):filedefault
REENTER DISPLAY FORMAT FOR ALL FILES (FILEDEFAULT):filedefault

L7 ANSWER 1 OF 9 HCAPLUS COPYRIGHT 2004 ACS on STN

L4 ANSWER 3 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 20 Sep 2004

L4 ANSWER 4 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 03 Feb 2004

L4 ANSWER 5 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 28 May 2003

L4 ANSWER 6 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 28 May 2003

L4 ANSWER 7 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 04 Jan 2002

L4 ANSWER 8 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 26 Dec 2001

=> d 14 1-8 1c

L4 ANSWER 1 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

L4 ANSWER 2 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

L4 ANSWER 3 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

L4 ANSWER 4 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
LC STN Files: CA, CAPLUS, USPATFULL

L4 ANSWER 5 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
LC STN Files: CHEMCATS

L4 ANSWER 6 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
LC STN Files: CHEMCATS

L4 ANSWER 7 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
LC STN Files: CHEMCATS

L4 ANSWER 8 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
LC STN Files: CHEMCATS

=> fil hcaplus uspatfull chemcats

COST IN U.S. DOLLARS SINCE FILE
TOTAL ENTRY

SESSION FULL ESTIMATED COST 294.47
294.68

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE
TOTAL ENTRY

SESSION CA SUBSCRIBER PRICE -1.32
1.32

AN 2004:41501 HCAPLUS Full-text

DN 140:87744

TI Affinity small molecules for the EPO receptor

IN Olsson, Lennart; Naranda, Tatjana

PA Receptor, Inc., USA

SO PCT Int. Appl., 85 pp.

CODEN: PIXXD2

DT Patent

LA English

FAN.CNT 1

PATENT NO. KIND DATE APPLICATION NO.

DATE -----

PI WO 2004005323 A2 20040115 WO 2003-US21394

20030703 WO 2004005323 A3 20040701

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA,

CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD,

GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC,

LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ,

OM, PH, PL, PT, RO, RU, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT,

TZ, UA, UG, UZ, VN, YU, ZA, ZM, ZW

RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM,

AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK,

EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI,

SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN,

TD, TG US 2004171541 A1 20040902 US 2003-613754

20030702 US 2004116346 A1 20040617 US 2003-612885

20030703 PRAI US 2002-393360P P 20020703

US 2002-393361P P 20020703

US 2002-394110P P 20020703

OS MARPAT 140:87744

L7 ANSWER 2 OF 9 USPATFULL on STN

AN 2004:221770 USPATFULL Full-text

TI Affinity small molecules for the EPO receptor

IN Olsson, Lennart; Orinda, CA, UNITED STATES

PI Naranda, Tatjana, Mountain View, CA, UNITED STATES

AI US 2004171541 A1 20040902

PRAI US 2003-613754 A1 20030702 (10)

US 2002-393361P 20020703 (60)

US 2002-393360P 20020703 (60)

US 2002-394110P 20020703 (60)

DT Utility

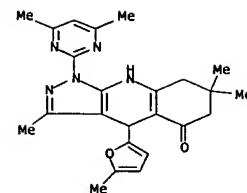
FS APPLICATION

LN.CNT 2046
 INCL INCLM: 514/012.000
 INCL: 514/247.000; 514/259.310; 514/292.000; 514/347.000;
 514/348.000
 NCL NCLM: 514/012.000
 NCLS: 514/247.000; 514/259.310; 514/292.000; 514/347.000;
 514/348.000
 IC [7]
 ICM: A61K038-18
 ICS: A61K031-50; A61K031-519; A61K031-44; A61K031-4745
 CAS INDEXING IS AVAILABLE FOR THIS PATENT.

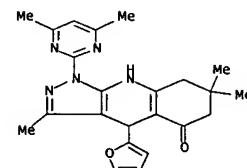
L7 ANSWER 3 OF 9 USPTFULL on STN
 AN 2004:152124 USPTFULL Full-text
 TI Affinity small molecules for the EPO receptor
 IN Olsson, Lennart, Orinda, CA, UNITED STATES
 Naranda, Tatjana, Mountain View, CA, UNITED STATES
 PI US 2004116346 A1 20040617
 AI US 2003-612885 A1 20030703 (10)
 PRAI US 2002-393361P 20020703 (60)
 US 2002-393360P 20020703 (60)
 US 2002-394110P 20020703 (60)
 DT Utility
 FS APPLICATION

LN.CNT 2000
 INCL INCLM: 514/012.000
 INCL: 514/247.000; 514/259.310; 514/292.000; 514/347.000;
 514/348.000
 NCL NCLM: 514/012.000
 NCLS: 514/247.000; 514/259.310; 514/292.000; 514/347.000;
 514/348.000
 IC [7]
 ICM: A61K038-17
 ICS: A61K031-519; A61K031-501; A61K031-4745; A61K031-44
 CAS INDEXING IS AVAILABLE FOR THIS PATENT.

L7 ANSWER 4 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2003:3262248 CHEMCATS
 Catalog Name (CO): Ambinter Screening Library
 Publication Date (PD): 1 Jan 2004
 Order Number (ON): T0509-4266
 Chemical Name (CN): 5H-Pyrazolo[3,4-b]quinolin-5-one,
 1-(4,6-dimethyl-2-pyrimidinyl)-1,4,6,7,8,9-
 hexahydro-
 3,7,7-trimethyl-4-(5-methyl-2-furanyl)-
 CAS Registry No. (RN): 521318-71-8
 Supplementary Term (ST): CHEMICAL LIBRARY
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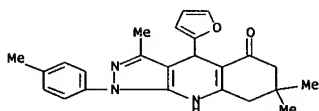


L7 ANSWER 5 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2003:3262236 CHEMCATS
 Catalog Name (CO): Ambinter Screening Library
 Publication Date (PD): 1 Jan 2004
 Order Number (ON): T0509-4092
 Chemical Name (CN): 5H-Pyrazolo[3,4-b]quinolin-5-one,
 1-(4,6-dimethyl-2-pyrimidinyl)-4-(2-
 furanyl)-
 1,4,6,7,8,9-hexahydro-3,7,7-trimethyl-
 CAS Registry No. (RN): 521284-01-5
 Supplementary Term (ST): CHEMICAL LIBRARY
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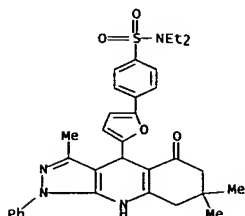


L7 ANSWER 6 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2003:2881171 CHEMCATS
 Catalog Name (CO): Enamine Screening Library
 Publication Date (PD): 30 Jun 2004
 Order Number (ON): T0508-1321
 Chemical Name (CN): 5H-Pyrazolo[3,4-b]quinolin-5-one,
 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-
 trimethyl-1-
 (4-methylphenyl)-
 CAS Registry No. (RN): 380450-98-6
 Supplementary Term (ST): CHEMICAL LIBRARY
 Structure :

Structure :

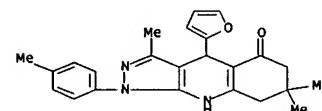


L7 ANSWER 7 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2003:2877484 CHEMCATS
 Catalog Name (CO): Enamine Screening Library
 Publication Date (PD): 30 Jun 2004
 Order Number (ON): T0505-5972
 Chemical Name (CN): Benzenesulfonamide, N,N-diethyl-4-[5-
 (4,5,6,7,8,9-
 hexahydro-3,7,7-trimethyl-5-oxo-1-phenyl)-
 pyrazolo[3,4-b]quinolin-4-yl]-2-furanyl]-
 CAS Registry No. (RN): 378189-53-8
 Supplementary Term (ST): CHEMICAL LIBRARY
 Structure :

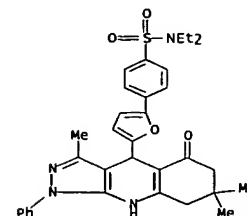


L7 ANSWER 8 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2002:1507742 CHEMCATS
 Catalog Name (CO): Ambinter Screening Library
 Publication Date (PD): 1 Jan 2004
 Order Number (ON): T0508-1321
 Chemical Name (CN): 5H-Pyrazolo[3,4-b]quinolin-5-one,
 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-
 trimethyl-1-
 (4-methylphenyl)-

CAS Registry No. (RN): 380450-98-6
 Supplementary Term (ST): CHEMICAL LIBRARY
 Structure :



L7 ANSWER 9 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2002:1491648 CHEMCATS
 Catalog Name (CO): Ambinter Screening Library
 Publication Date (PD): 1 Jan 2004
 Order Number (ON): T0505-5972
 Chemical Name (CN): Benzenesulfonamide, N,N-diethyl-4-[5-
 (4,5,6,7,8,9-
 hexahydro-3,7,7-trimethyl-5-oxo-1-phenyl)-
 pyrazolo[3,4-b]quinolin-4-yl]-2-furanyl]-
 CAS Registry No. (RN): 378189-53-8
 Supplementary Term (ST): CHEMICAL LIBRARY
 Structure :



=> d 17 1-9 pd
 NO VALID FORMATS ENTERED FOR FILE 'HCAPLUS'
 NO VALID FORMATS ENTERED FOR FILE 'USPTFULL'
 In a multife environment, each file must have at least one valid
 format requested. Refer to file specific help messages or the
 STNGUIDE file for information on formats available in individual


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files.
REENTER DISPLAY FORMAT FOR ALL FILES (FILEDEFAULT):end

=> index biosci
FILE 'DRUGMONOG' ACCESS NOT AUTHORIZED
COST IN U.S. DOLLARS          SINCE FILE
TOTAL                          ENTRY
SESSION
FULL ESTIMATED COST          19.21
313.89

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE
TOTAL                          ENTRY
SESSION
CA SUBSCRIBER PRICE          0.00 -
1.32

INDEX 'ADISCTI, ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, ANTE,
AQUALINE,
AQUASCI, BIOBUSINESS, BIOCERAMICS, BIOENG, BIOSIS, BIOTECHABS,
BIOTECHDS,
BIOTECHNO, CABA, CANCERLIT, CAPLUS, CEABA-VTB, CEN, CIN,
CONFSCI, CROPB,
CROPU, DDFB, DDFU, DGENE, DISSABS, ...' ENTERED AT 10:38:10 ON
28 DEC 2004

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75 FILES IN THE FILE LIST IN STINDEX

Enter SET DETAIL ON to see search term postings or to view search error messages that display as 0* with SET DETAIL OFF.

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=> s (epo or (erythropoietin receptor))
FILE 'ADISCTI'
  68 EPO
  3 EPOS
  71 EPO
    (EPO OR EPOS)
  641 ERYTHROPOIETIN
  1 ERYTHROPOIETINS
  642 ERYTHROPOIETIN
    (ERYTHROPOIETIN OR ERYTHROPOIETINS)
  77336 RECEPTOR
  5437 RECEPTORS
  79811 RECEPTOR
    (RECEPTOR OR RECEPTORS)
  3 ERYTHROPOIETIN RECEPTOR
    (ERYTHROPOIETIN(W)RECEPTOR)
  73 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'ADISINSIGHT'
  24 EPO
  2 EPOS
  25 EPO
    (EPO OR EPOS)
  57 "ERYTHROPOIETIN"
  15 "ERYTHROPOIETINS"
    (RECEPTOR OR RECEPTORS)
  0 ERYTHROPOIETIN RECEPTOR
    (ERYTHROPOIETIN(W)RECEPTOR)
  22 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'AQUALINE'
  1 EPO
  2 EPOS
  3 EPO
    (EPO OR EPOS)
  1 ERYTHROPOIETIN
  241 RECEPTOR
  286 RECEPTORS
  395 RECEPTOR
    (RECEPTOR OR RECEPTORS)
  0 ERYTHROPOIETIN RECEPTOR
    (ERYTHROPOIETIN(W)RECEPTOR)
  3 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'AQUASCI'
  36 EPO
  63 EPOS
  99 EPO
    (EPO OR EPOS)
  10 "ERYTHROPOIETIN"
  5274 "RECEPTOR"
  4864 "RECEPTORS"
  7710 "RECEPTOR"
    ("RECEPTOR" OR "RECEPTORS")
  0 ERYTHROPOIETIN RECEPTOR
    ("ERYTHROPOIETIN" (W) "RECEPTOR")
  99 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'BIOBUSINESS'
  387 EPO
  18 EPOS
  404 EPO
    (EPO OR EPOS)
  588 "ERYTHROPOIETIN"
  2 "ERYTHROPOIETINS"
  590 "ERYTHROPOIETIN"
    ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
  6807 "RECEPTOR"
  2441 "RECEPTORS"
  7936 "RECEPTOR"
    ("RECEPTOR" OR "RECEPTORS")
  7 ERYTHROPOIETIN RECEPTOR
    ("ERYTHROPOIETIN" (W) "RECEPTOR")
  410 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'BIOCERAMICS'
  1192 EPO
  5 EPOS
  1197 EPO
    (EPO OR EPOS)
  837 ERYTHROPOIETIN
  1 ERYTHROPOIETINS
  837 ERYTHROPOIETIN
    (ERYTHROPOIETIN OR ERYTHROPOIETINS)
  3060 RECEPTOR
  1032 RECEPTORS

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  57 "ERYTHROPOIETIN"
    ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
  4128 "RECEPTOR"
  1761 "RECEPTORS"
  4546 "RECEPTOR"
    ("RECEPTOR" OR "RECEPTORS")
  6 ERYTHROPOIETIN RECEPTOR
    ("ERYTHROPOIETIN" (W) "RECEPTOR")
  27 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'ADISNEWS'
  16 EPO
  162 ERYTHROPOIETIN
  5 ERYTHROPOIETINS
  163 ERYTHROPOIETIN
    (ERYTHROPOIETIN OR ERYTHROPOIETINS)
  1962 RECEPTOR
  556 RECEPTORS
  2258 RECEPTOR
    (RECEPTOR OR RECEPTORS)
  3 ERYTHROPOIETIN RECEPTOR
    (ERYTHROPOIETIN(W)RECEPTOR)
  19 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'AGRICOLA'
  35 EPO
  26 EPOS
  60 EPO
    (EPO OR EPOS)
  136 ERYTHROPOIETIN
  8600 RECEPTOR
  9839 RECEPTORS
  13992 RECEPTOR
    (RECEPTOR OR RECEPTORS)
  3 ERYTHROPOIETIN RECEPTOR
    (ERYTHROPOIETIN(W)RECEPTOR)
  63 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'ANABSTR'
  26 EPO
  27 EPOS
  52 EPO
    (EPO OR EPOS)
  72 ERYTHROPOIETIN
  1181 RECEPTOR
  381 RECEPTORS
  1391 RECEPTOR
    (RECEPTOR OR RECEPTORS)
  1 ERYTHROPOIETIN RECEPTOR
    (ERYTHROPOIETIN(W)RECEPTOR)
  53 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'ANTE'
  10 EPO
  12 EPOS
  22 EPO
    (EPO OR EPOS)
  7 ERYTHROPOIETIN
  91 RECEPTOR
  73 RECEPTORS
  150 RECEPTOR
    (RECEPTOR OR RECEPTORS)
  3917 RECEPTOR
    (RECEPTOR OR RECEPTORS)
  0 ERYTHROPOIETIN RECEPTOR
    (ERYTHROPOIETIN(W)RECEPTOR)
  1197 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'BIOENG'
  180 EPO
  7 EPOS
  183 EPO
    (EPO OR EPOS)
  335 ERYTHROPOIETIN
  9 ERYTHROPOIETINS
  336 ERYTHROPOIETIN
    (ERYTHROPOIETIN OR ERYTHROPOIETINS)
  11378 RECEPTOR
  8432 RECEPTORS
  14157 RECEPTOR
    (RECEPTOR OR RECEPTORS)
  28 ERYTHROPOIETIN RECEPTOR
    (ERYTHROPOIETIN(W)RECEPTOR)
  199 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'BIOSIS'
  5622 EPO
  222 EPOS
  5822 EPO
    (EPO OR EPOS)
  17873 ERYTHROPOIETIN
  64 ERYTHROPOIETINS
  17897 ERYTHROPOIETIN
    (ERYTHROPOIETIN OR ERYTHROPOIETINS)
  651542 RECEPTOR
  323743 RECEPTORS
  780807 RECEPTOR
    (RECEPTOR OR RECEPTORS)
  1356 ERYTHROPOIETIN RECEPTOR
    (ERYTHROPOIETIN(W)RECEPTOR)
  6546 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'BIOTECHABS'
  356 EPO
  8 EPOS
  356 EPO
    (EPO OR EPOS)
  981 ERYTHROPOIETIN
  6 ERYTHROPOIETINS
  983 ERYTHROPOIETIN
    (ERYTHROPOIETIN OR ERYTHROPOIETINS)
  16255 RECEPTOR
  4183 RECEPTORS
  17462 RECEPTOR
    (RECEPTOR OR RECEPTORS)
  85 ERYTHROPOIETIN RECEPTOR
    (ERYTHROPOIETIN(W)RECEPTOR)
  407 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'BIOTECHDS'
  356 EPO
  8 EPOS
  356 EPO

```

(EPO OR EPOS)
 981 ERYTHROPOIETIN
 6 ERYTHROPOIETINS
 983 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 16255 RECEPTOR
 4183 RECEPTORS
 17462 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 85 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 407 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'BIOTECHNO'
 2598 EPO
 33 EPOS
 2614 EPO
 (EPO OR EPOS)
 9222 ERYTHROPOIETIN
 29 ERYTHROPOIETINS
 9222 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 202158 RECEPTOR
 78475 RECEPTORS
 213334 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 774 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 2960 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CABA'
 247 EPO
 47 EPOS
 294 EPO
 (EPO OR EPOS)
 539 ERYTHROPOIETIN
 30692 RECEPTOR
 27815 RECEPTORS
 42614 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 21 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 308 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CANCERLIT'
 3530 EPO
 31 EPOS
 3544 EPO
 (EPO OR EPOS)
 11104 ERYTHROPOIETIN
 30 ERYTHROPOIETINS
 11104 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 122617 RECEPTOR
 117416 RECEPTORS
 160576 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 501 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 3775 (EPO OR (ERYTHROPOIETIN RECEPTOR))

539 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CONFSCI'
 87 EPO
 10 EPOS
 97 EPO
 (EPO OR EPOS)
 525 "ERYTHROPOIETIN"
 2 "ERYTHROPOIETINS"
 527 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 17313 "RECEPTOR"
 8593 "RECEPTORS"
 25653 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 71 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN"(W)"RECEPTOR")
 165 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CROBP'
 0 EPO
 9 EPOS
 9 EPO
 (EPO OR EPOS)
 0 ERYTHROPOIETIN
 160 RECEPTOR
 35 RECEPTORS
 175 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 9 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CROPU'
 55 EPO
 12 EPOS
 67 EPO
 (EPO OR EPOS)
 1 ERYTHROPOIETIN
 1368 RECEPTOR
 570 RECEPTORS
 1566 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 67 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'DDFB'
 2 EPO
 843 ERYTHROPOIETIN
 15280 RECEPTOR
 4633 RECEPTORS
 16551 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 3 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'DDFU'
 1419 EPO
 3 EPOS
 1421 EPO

FILE 'CAPLUS'
 5521 EPO
 131 EPOS
 5625 EPO
 (EPO OR EPOS)
 11125 ERYTHROPOIETIN
 520 ERYTHROPOIETINS
 11155 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 582865 RECEPTOR
 534615 RECEPTORS
 693859 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1237 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 6216 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CEABA-VTB'
 106 EPO
 16 EPOS
 122 EPO
 (EPO OR EPOS)
 191 ERYTHROPOIETIN
 9 ERYTHROPOIETINS
 196 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 2128 RECEPTOR
 850 RECEPTORS
 2456 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 4 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 124 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CEN'
 33 EPO
 67 "ERYTHROPOIETIN"
 439 "RECEPTOR"
 381 "RECEPTORS"
 630 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 1 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN"(W)"RECEPTOR")
 34 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CIN'
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 6 EPOS
 536 EPO
 (EPO OR EPOS)
 544 "ERYTHROPOIETIN"
 4 "ERYTHROPOIETINS"
 547 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 4037 "RECEPTOR"
 1359 "RECEPTORS"
 4824 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 7 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN"(W)"RECEPTOR")

(EPO OR EPOS)
 3133 ERYTHROPOIETIN
 9 ERYTHROPOIETINS
 3133 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 99698 RECEPTOR
 44104 RECEPTORS
 111042 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 56 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 1441 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'DGENE'
 15416 EPO
 68 EPOS
 15418 EPO
 (EPO OR EPOS)
 18662 ERYTHROPOIETIN
 21 ERYTHROPOIETINS
 18680 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 496912 RECEPTOR
 170378 RECEPTORS
 537954 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1551 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 15707 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'DISSABS'
 152 EPO
 51 EPOS
 200 EPO
 (EPO OR EPOS)
 240 ERYTHROPOIETIN
 22949 RECEPTOR
 13830 RECEPTORS
 27515 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 35 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 216 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'DRUGB'
 2 EPO
 843 ERYTHROPOIETIN
 15280 RECEPTOR
 4633 RECEPTORS
 16551 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 3 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'DRUGMONOG2'
 74 EPO
 2 EPOS
 76 EPO
 (EPO OR EPOS)
 1 ERYTHROPOIETIN

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0 RECEPTOR
0 ERYTHROPOIETIN RECEPTOR
  (ERYTHROPOIETIN(W)RECEPTOR)
76 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'DRUGU'
1633 EPO
7 EPOS
1637 EPO
  (EPO OR EPOS)
3396 ERYTHROPOIETIN
10 ERYTHROPOIETINS
3396 ERYTHROPOIETIN
  (ERYTHROPOIETIN OR ERYTHROPOIETINS)
110636 RECEPTOR
54912 RECEPTORS
125028 RECEPTOR
  (RECEPTOR OR RECEPTORS)
64 ERYTHROPOIETIN RECEPTOR
  (ERYTHROPOIETIN(W)RECEPTOR)
1657 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'EMBAL'
52 EPO
2 EPOS
54 EPO
  (EPO OR EPOS)
106 ERYTHROPOIETIN
2 ERYTHROPOIETINS
108 ERYTHROPOIETIN
  (ERYTHROPOIETIN OR ERYTHROPOIETINS)
5008 RECEPTOR
2397 RECEPTORS
5899 RECEPTOR
  (RECEPTOR OR RECEPTORS)
6 ERYTHROPOIETIN RECEPTOR
  (ERYTHROPOIETIN(W)RECEPTOR)
57 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'EMBASE'
4482 EPO
111 EPOS
4572 EPO
  (EPO OR EPOS)
18291 "ERYTHROPOIETIN"
51 "ERYTHROPOIETINS"
18292 "ERYTHROPOIETIN"
  ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
714332 "RECEPTOR"
260226 "RECEPTORS"
755991 "RECEPTOR"
  ("RECEPTOR" OR "RECEPTORS")
1098 ERYTHROPOIETIN RECEPTOR
  ("ERYTHROPOIETIN"(W)"RECEPTOR")
5095 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'ESBIOBASE'
1925 EPO
44 EPOS
1960 EPO
  (EPO OR EPOS)

FILE 'FSTA'
21 EPO
5 EPOS
25 EPO
  (EPO OR EPOS)
1 ERYTHROPOIETIN
1012 RECEPTOR
434 RECEPTORS
1297 RECEPTOR
  (RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
  (ERYTHROPOIETIN(W)RECEPTOR)
25 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'GENBANK'
6183 EPO
6652 "ERYTHROPOIETIN"
201129 "RECEPTOR"
289 ERYTHROPOIETIN RECEPTOR
  ("ERYTHROPOIETIN"(W)"RECEPTOR")
6457 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'HEALSAFE'
7 EPO
19 "ERYTHROPOIETIN"
460 "RECEPTOR"
311 "RECEPTORS"
657 "RECEPTOR"
  ("RECEPTOR" OR "RECEPTORS")
0 ERYTHROPOIETIN RECEPTOR
  ("ERYTHROPOIETIN"(W)"RECEPTOR")
7 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'IFIPAT'
716 EPO
11 EPOS
725 EPO
  (EPO OR EPOS)
1216 ERYTHROPOIETIN
24 ERYTHROPOIETINS
1235 ERYTHROPOIETIN
  (ERYTHROPOIETIN OR ERYTHROPOIETINS)
32172 RECEPTOR
11668 RECEPTORS
36258 RECEPTOR
  (RECEPTOR OR RECEPTORS)
68 ERYTHROPOIETIN RECEPTOR
  (ERYTHROPOIETIN(W)RECEPTOR)
758 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'IMSDRUGNEWS'
70 EPO
161 "ERYTHROPOIETIN"
3340 "RECEPTOR"
836 "RECEPTORS"
3873 "RECEPTOR"
  ("RECEPTOR" OR "RECEPTORS")
7 ERYTHROPOIETIN RECEPTOR
  ("ERYTHROPOIETIN"(W)"RECEPTOR")
74 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'IMSPRODUCT'

3664 ERYTHROPOIETIN
11 ERYTHROPOIETINS
3666 ERYTHROPOIETIN
  (ERYTHROPOIETIN OR ERYTHROPOIETINS)
239241 RECEPTOR
129772 RECEPTORS
280062 RECEPTOR
  (RECEPTOR OR RECEPTORS)
433 ERYTHROPOIETIN RECEPTOR
  (ERYTHROPOIETIN(W)RECEPTOR)
2150 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'FEDRIP'
118 EPO
2 EPOS
120 EPO
  (EPO OR EPOS)
243 ERYTHROPOIETIN
16896 RECEPTOR
9138 RECEPTORS
19085 RECEPTOR
  (RECEPTOR OR RECEPTORS)
30 ERYTHROPOIETIN RECEPTOR
  (ERYTHROPOIETIN(W)RECEPTOR)
139 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'FOMAD'
1 EPO
12 EPOS
13 EPO
  (EPO OR EPOS)
1 ERYTHROPOIETIN
2 RECEPTOR
1 RECEPTORS
2 RECEPTOR
  (RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
  (ERYTHROPOIETIN(W)RECEPTOR)
13 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'FOREGE'
0 EPO
0 ERYTHROPOIETIN
0 RECEPTOR
0 ERYTHROPOIETIN RECEPTOR
  (ERYTHROPOIETIN(W)RECEPTOR)
0 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'FROSTI'
15 EPO
15 EPOS
29 EPO
  (EPO OR EPOS)
9 ERYTHROPOIETIN
1085 RECEPTOR
784 RECEPTORS
1582 RECEPTOR
  (RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
  (ERYTHROPOIETIN(W)RECEPTOR)
29 (EPO OR (ERYTHROPOIETIN RECEPTOR))

31 EPO
2 EPOS
33 EPO
  (EPO OR EPOS)
219 "ERYTHROPOIETIN"
44 "RECEPTOR"
8 "RECEPTORS"
52 "RECEPTOR"
  ("RECEPTOR" OR "RECEPTORS")
0 ERYTHROPOIETIN RECEPTOR
  ("ERYTHROPOIETIN"(W)"RECEPTOR")
33 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'IMSRESEARCH'
32 EPO
70 "ERYTHROPOIETIN"
2213 "RECEPTOR"
1048 "RECEPTORS"
2816 "RECEPTOR"
  ("RECEPTOR" OR "RECEPTORS")
6 ERYTHROPOIETIN RECEPTOR
  ("ERYTHROPOIETIN"(W)"RECEPTOR")
36 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'JICST-EPLUS'
621 EPO
17 EPOS
635 EPO
  (EPO OR EPOS)
3631 ERYTHROPOIETIN
4 ERYTHROPOIETINS
3631 ERYTHROPOIETIN
  (ERYTHROPOIETIN OR ERYTHROPOIETINS)
73555 RECEPTOR
12647 RECEPTORS
76150 RECEPTOR
  (RECEPTOR OR RECEPTORS)
192 ERYTHROPOIETIN RECEPTOR
  (ERYTHROPOIETIN(W)RECEPTOR)
800 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'KOSMET'
5 EPO
1 EPOS
6 EPO
  (EPO OR EPOS)
8 ERYTHROPOIETIN
490 RECEPTOR
555 RECEPTORS
710 RECEPTOR
  (RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
  (ERYTHROPOIETIN(W)RECEPTOR)
6 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'LIFESCI'
863 EPO
50 EPOS
907 EPO
  (EPO OR EPOS)
1458 "ERYTHROPOIETIN"

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11 "ERYTHROPOIETINS"
1458 "ERYTHROPOIETIN"
    ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
167986 "RECEPTOR"
153110 "RECEPTORS"
216519 "RECEPTOR"
    ("RECEPTOR" OR "RECEPTORS")
281 ERYTHROPOIETIN RECEPTOR
    ("ERYTHROPOIETIN"(W)"RECEPTOR")
1047 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'MEDICINF'
    4 EPO
    7 EPOS
    11 EPO
        (EPO OR EPOS)
    22 ERYTHROPOIETIN
    1 ERYTHROPOIETINS
    23 ERYTHROPOIETIN
        (ERYTHROPOIETIN OR ERYTHROPOIETINS)
    324 RECEPTOR
    571 RECEPTORS
    827 RECEPTOR
        (RECEPTOR OR RECEPTORS)
    0 ERYTHROPOIETIN RECEPTOR
        (ERYTHROPOIETIN(W)RECEPTOR)
    11 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'MEDLINE'
    4856 EPO
    99 EPOS
    4929 EPO
        (EPO OR EPOS)
    16393 ERYTHROPOIETIN
    53 ERYTHROPOIETINS
    16397 ERYTHROPOIETIN
        (ERYTHROPOIETIN OR ERYTHROPOIETINS)
    479689 RECEPTOR
    514892 RECEPTORS
    673211 RECEPTOR
        (RECEPTOR OR RECEPTORS)
    714 ERYTHROPOIETIN RECEPTOR
        (ERYTHROPOIETIN(W)RECEPTOR)
    5281 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'NIOSHITIC'
    15 EPO
    89 ERYTHROPOIETIN
    1 ERYTHROPOIETINS
    89 ERYTHROPOIETIN
        (ERYTHROPOIETIN OR ERYTHROPOIETINS)
    1641 RECEPTOR
    1030 RECEPTORS
    2222 RECEPTOR
        (RECEPTOR OR RECEPTORS)
    0 ERYTHROPOIETIN RECEPTOR
        (ERYTHROPOIETIN(W)RECEPTOR)
    15 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'NTIS'
    46 EPO

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64 EPOS
108 EPO
    (EPO OR EPOS)
121 ERYTHROPOIETIN
25 ERYTHROPOIETINS
122 ERYTHROPOIETIN
    (ERYTHROPOIETIN OR ERYTHROPOIETINS)
4927 RECEPTOR
3307 RECEPTORS
6501 RECEPTOR
    (RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
    (ERYTHROPOIETIN(W)RECEPTOR)
108 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'NUTRACEUT'
    11 EPO
    0 ERYTHROPOIETIN
    10 RECEPTOR
    13 RECEPTORS
    22 RECEPTOR
        (RECEPTOR OR RECEPTORS)
    0 ERYTHROPOIETIN RECEPTOR
        (ERYTHROPOIETIN(W)RECEPTOR)
    11 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'OCEAN'
    24 EPO
    51 EPOS
    75 EPO
        (EPO OR EPOS)
    1 "ERYTHROPOIETIN"
    601 "RECEPTOR"
    581 "RECEPTORS"
    950 "RECEPTOR"
        ("RECEPTOR" OR "RECEPTORS")
    0 ERYTHROPOIETIN RECEPTOR
        ("ERYTHROPOIETIN"(W)"RECEPTOR")
    75 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'PASCAL'
    2055 EPO
    127 EPOS
    2173 EPO
        (EPO OR EPOS)
    6424 ERYTHROPOIETIN
    22 ERYTHROPOIETINS
    6427 ERYTHROPOIETIN
        (ERYTHROPOIETIN OR ERYTHROPOIETINS)
    304437 RECEPTOR
    102112 RECEPTORS
    327238 RECEPTOR
        (RECEPTOR OR RECEPTORS)
    334 ERYTHROPOIETIN RECEPTOR
        (ERYTHROPOIETIN(W)RECEPTOR)
    2342 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'PCTGEN'
    0 EPO
    0 ERYTHROPOIETIN
    8549 RECEPTOR

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694 RECEPTORS
9243 RECEPTOR
    (RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
    (ERYTHROPOIETIN(W)RECEPTOR)
0 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'PHAR'
    72 EPO
    85 "ERYTHROPOIETIN"
    9931 "RECEPTOR"
    1111 "RECEPTORS"
    10104 "RECEPTOR"
        ("RECEPTOR" OR "RECEPTORS")
    47 ERYTHROPOIETIN RECEPTOR
        ("ERYTHROPOIETIN"(W)"RECEPTOR")
    73 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'PHARMAML'
    99 EPO
    3 EPOS
    101 EPO
        (EPO OR EPOS)
    202 ERYTHROPOIETIN
    5 ERYTHROPOIETINS
    206 ERYTHROPOIETIN
        (ERYTHROPOIETIN OR ERYTHROPOIETINS)
    1768 RECEPTOR
    465 RECEPTORS
    2031 RECEPTOR
        (RECEPTOR OR RECEPTORS)
    0 ERYTHROPOIETIN RECEPTOR
        (ERYTHROPOIETIN(W)RECEPTOR)
    101 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'PHIC'
    7 EPO
    6 "ERYTHROPOIETIN"
    18 "RECEPTOR"
    5 "RECEPTORS"
    23 "RECEPTOR"
        ("RECEPTOR" OR "RECEPTORS")
    0 ERYTHROPOIETIN RECEPTOR
        ("ERYTHROPOIETIN"(W)"RECEPTOR")
    7 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'PHIN'
    841 EPO
    22 EPOS
    859 EPO
        (EPO OR EPOS)
    1142 "ERYTHROPOIETIN"
    18 "ERYTHROPOIETINS"
    1153 "ERYTHROPOIETIN"
        ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
    4441 "RECEPTOR"
    2186 "RECEPTORS"
    5782 "RECEPTOR"
        ("RECEPTOR" OR "RECEPTORS")
    3 ERYTHROPOIETIN RECEPTOR
        ("ERYTHROPOIETIN"(W)"RECEPTOR")

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861 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'PROMT'
    2713 EPO
    1902 EPOS
    4590 EPO
        (EPO OR EPOS)
    2025 "ERYTHROPOIETIN"
    34 "ERYTHROPOIETINS"
    2050 "ERYTHROPOIETIN"
        ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
    17544 "RECEPTOR"
    9487 "RECEPTORS"
    22972 "RECEPTOR"
        ("RECEPTOR" OR "RECEPTORS")
    27 ERYTHROPOIETIN RECEPTOR
        ("ERYTHROPOIETIN"(W)"RECEPTOR")
    4612 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'PROUDDDR'
    22 EPO
    43 "ERYTHROPOIETIN"
    2 "ERYTHROPOIETINS"
    44 "ERYTHROPOIETIN"
        ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
    36740 "RECEPTOR"
    19875 "RECEPTORS"
    44173 "RECEPTOR"
        ("RECEPTOR" OR "RECEPTORS")
    4 ERYTHROPOIETIN RECEPTOR
        ("ERYTHROPOIETIN"(W)"RECEPTOR")
    22 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'PS'
    0 EPO
    1 EPOS
    1 EPO
        (EPO OR EPOS)
    0 ERYTHROPOIETIN
    29 RECEPTOR
    1 RECEPTORS
    30 RECEPTOR
        (RECEPTOR OR RECEPTORS)
    0 ERYTHROPOIETIN RECEPTOR
        (ERYTHROPOIETIN(W)RECEPTOR)
    1 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'RDISCLOSURE'
    21 EPO
    2 EPOS
    23 EPO
        (EPO OR EPOS)
    1 ERYTHROPOIETIN
    93 RECEPTOR
    20 RECEPTORS
    106 RECEPTOR
        (RECEPTOR OR RECEPTORS)
    0 ERYTHROPOIETIN RECEPTOR
        (ERYTHROPOIETIN(W)RECEPTOR)
    23 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'SCISEARCH'

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4531 EPO
222 EPOS
4734 EPO
      (EPO OR EPOS)
16279 ERYTHROPOIETIN
46 ERYTHROPOIETINS
16295 ERYTHROPOIETIN
      (ERYTHROPOIETIN OR ERYTHROPOIETINS)
583687 RECEPTOR
306298 RECEPTORS
721251 RECEPTOR
      (RECEPTOR OR RECEPTORS)
1592 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
5851 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'SYNTHLINE'
1 EPO
0 "ERYTHROPOIETIN"
798 "RECEPTOR"
103 "RECEPTORS"
876 "RECEPTOR"
      ("RECEPTOR" OR "RECEPTORS")
0 ERYTHROPOIETIN RECEPTOR
      ("ERYTHROPOIETIN"(W)"RECEPTOR")
1 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'TOXCENTER'
2126 EPO
34 EPOS
2145 EPO
      (EPO OR EPOS)
6597 ERYTHROPOIETIN
72 ERYTHROPOIETINS
6605 ERYTHROPOIETIN
      (ERYTHROPOIETIN OR ERYTHROPOIETINS)
264727 RECEPTOR
166955 RECEPTORS
322116 RECEPTOR
      (RECEPTOR OR RECEPTORS)
285 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
2274 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'USPATFULL'
13847 EPO
184 EPOS
13982 EPO
      (EPO OR EPOS)
7989 ERYTHROPOIETIN
1884 ERYTHROPOIETINS
9673 ERYTHROPOIETIN
      (ERYTHROPOIETIN OR ERYTHROPOIETINS)
104220 RECEPTOR
75854 RECEPTORS
118925 RECEPTOR
      (RECEPTOR OR RECEPTORS)
773 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
14435 (EPO OR (ERYTHROPOIETIN RECEPTOR))

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      (ERYTHROPOIETIN OR ERYTHROPOIETINS)
41921 RECEPTOR
13906 RECEPTORS
46849 RECEPTOR
      (RECEPTOR OR RECEPTORS)
80 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
659 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'WPIFV'
10 EPO
15 ERYTHROPOIETIN
354 RECEPTOR
89 RECEPTORS
389 RECEPTOR
      (RECEPTOR OR RECEPTORS)
2 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
10 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'WPIINDEX'
580 EPO
36 EPOS
612 EPO
      (EPO OR EPOS)
1446 ERYTHROPOIETIN
16 ERYTHROPOIETINS
1458 ERYTHROPOIETIN
      (ERYTHROPOIETIN OR ERYTHROPOIETINS)
41921 RECEPTOR
13906 RECEPTORS
46849 RECEPTOR
      (RECEPTOR OR RECEPTORS)
80 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
659 (EPO OR (ERYTHROPOIETIN RECEPTOR))

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L8 QUE (EPO OR (ERYTHROPOIETIN RECEPTOR))

=> DIS HIST

(FILE 'HOME' ENTERED AT 10:25:10 ON 28 DEC 2004)

FILE 'REGISTRY' ENTERED AT 10:25:22 ON 28 DEC 2004

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L1 STRUCTURE UPLOADED
L2 0 S L1 SAM
L3 0 S L1 FAM
L4 8 S L1 FUL
L5 1 S L1 FUL FAM
L6 0 L5 NOT L4

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FILE 'HCAPLUS, USPATFULL, CHEMCATS' ENTERED AT 10:35:54 ON 28 DEC 2004

L7 9 S L4

INDEX 'ADISCTI, ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, ANTE, AQUALINE, AQUASCI, BIOBUSINESS, BIOCOMMERCE, BIOENG, BIOSIS, BIOTECHABS, BIOTECHDS,

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FILE 'USPAT2'
807 EPO
13 EPOS
818 EPO
      (EPO OR EPOS)
477 ERYTHROPOIETIN
30 ERYTHROPOIETINS
494 ERYTHROPOIETIN
      (ERYTHROPOIETIN OR ERYTHROPOIETINS)
7293 RECEPTOR
5119 RECEPTORS
8267 RECEPTOR
      (RECEPTOR OR RECEPTORS)
32 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
836 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'VETB'
0 EPO
1 ERYTHROPOIETIN
77 RECEPTOR
25 RECEPTORS
82 RECEPTOR
      (RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
0 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'VETU'
23 EPO
65 ERYTHROPOIETIN
970 RECEPTOR
641 RECEPTORS
1302 RECEPTOR
      (RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
23 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'WATER'
0 EPO
2 EPOS
2 EPO
      (EPO OR EPOS)
1 ERYTHROPOIETIN
471 RECEPTOR
265 RECEPTORS
641 RECEPTOR
      (RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
2 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'WPIDS'
580 EPO
36 EPOS
612 EPO
      (EPO OR EPOS)
1446 ERYTHROPOIETIN
16 ERYTHROPOIETINS
1458 ERYTHROPOIETIN

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BIOTECHNO, CABA, CANCERLIT, CAPLUS, CEABA-VTB, CEN, CIN, CONFSCI, CROPB, CROPU, DDFB, DDFU, DGENE, DISSABS, ...' ENTERED AT 10:38:10 ON 28 DEC 2004

SEA (EPO OR (ERYTHROPOIETIN RECEPTOR))

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73 FILE ADISCTI
27 FILE ADISINSIGHT
19 FILE ADISNEWS
63 FILE AGRICOLA
53 FILE ANABSTR
22 FILE ANTE
3 FILE AQUALINE
99 FILE AQUASCI
410 FILE BIOBUSINESS
1197 FILE BIOCOMMERCE
199 FILE BIOENG
6546 FILE BIOSIS
407 FILE BIOTECHABS
407 FILE BIOTECHDS
2960 FILE BIOTECHNO
308 FILE CABA
3775 FILE CANCERLIT
6216 FILE CAPLUS
124 FILE CEABA-VTB
34 FILE CEN
539 FILE CIN
165 FILE CONFSCI
9 FILE CROPB
67 FILE CROPU
3 FILE DDFB
1441 FILE DDFU
15707 FILE DGENE
216 FILE DISSABS
3 FILE DRUGB
76 FILE DRUGMONOG2
1657 FILE DRUGU
57 FILE EMBAL
5095 FILE EMBASE
2150 FILE ESBIODBASE
139 FILE FEDRIP
13 FILE FOMAD
29 FILE FROSTI
25 FILE FSTA
6457 FILE GENBANK
7 FILE HEALSAFE
758 FILE IFIPAT
74 FILE IMSDRUGNEWS
33 FILE IMSPRODUCT
36 FILE IMSRESEARCH
800 FILE JICST-EPLUS
6 FILE KOSMET
1047 FILE LIFESCI
11 FILE MEDICONF
5281 FILE MEDLINE
15 FILE NIOSHTIC
108 FILE NTIS

```

```

11 FILE NUTRACEUT
75 FILE OCEAN
2342 FILE PASCAL
73 FILE PHAR
101 FILE PHARMAML
7 FILE PHIC
861 FILE PHIN
4612 FILE PROMT
22 FILE PROUSDDR
1 FILE PS
23 FILE RDISCLOSURE
5851 FILE SCISEARCH
1 FILE SYNTHLINE
2274 FILE TOXCENTER
14435 FILE USPATFULL
836 FILE USPAT2
23 FILE VETU
2 FILE WATER
659 FILE WPIDS
10 FILE WPIFV
659 FILE WPINDE
L8 QUE (EPO OR (ERYTHROPOIETIN RECEPTOR))
-----

```

=>

---Logging off of STN---

=>

Executing the logoff script...

=> LOG Y

COST IN U.S. DOLLARS	SINCE FILE	
TOTAL	ENTRY	
SESSION		
FULL ESTIMATED COST	11.97	
325.86		
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	
TOTAL	ENTRY	
SESSION		
CA SUBSCRIBER PRICE	0.00	-
1.32		

STN INTERNATIONAL LOGOFF AT 10:50:37 ON 28 DEC 2004